

## Connecting via Winsock to STN

Trying 3106016892...Open

Welcome to STN International! Enter x:x

LOGINID: ssspta1623paz

PASSWORD :

TERMINAL (ENTER 1, 2, 3, OR ?):2

NEWS 1 Web Page URLs for STN Seminar Schedule - N. America  
NEWS 2 Sep 17 IMSworld Pharmaceutical Company Directory name change to PHARMASEARCH  
NEWS 3 Oct 09 Korean abstracts now included in Derwent World Patents Index  
NEWS 4 Oct 09 Number of Derwent World Patents Index updates increased  
NEWS 5 Oct 15 Calculated properties now in the REGISTRY/ZREGISTRY File  
NEWS 6 Oct 22 Over 1 million reactions added to CASREACT  
NEWS 7 Oct 22 DGENE GETSIM has been improved  
NEWS 8 Oct 29 AAASD no longer available  
NEWS 9 Nov 19 New Search Capabilities USPATFULL and USPAT2  
NEWS 10 Nov 19 TOXCENTER(SM) - new toxicology file now available on STN  
NEWS 11 Nov 29 COPPERLIT now available on STN  
NEWS 12 Nov 29 DWPI revisions to NTIS and US Provisional Numbers  
NEWS 13 Nov 30 Files VETU and VETB to have open access  
NEWS 14 Dec 10 WPINDEX/WPIDS/WPIX New and Revised Manual Codes for 2002  
NEWS 15 Dec 10 DGENE BLAST Homology Search  
NEWS 16 Dec 17 WELDASEARCH now available on STN  
NEWS 17 Dec 17 STANDARDS now available on STN  
NEWS 18 Dec 17 New fields for DPCI  
NEWS 19 Dec 19 CAS Roles modified  
NEWS 20 Dec 19 1907-1946 data and page images added to CA and CAplus  
NEWS 21 Jan 25 BLAST(R) searching in REGISTRY available in STN on the Web  
NEWS 22 Jan 25 Searching with the P indicator for Preparations  
NEWS 23 Jan 29 FSTA has been reloaded and moves to weekly updates  
NEWS 24 Feb 01 DKILIT now produced by FIZ Karlsruhe and has a new update frequency  
NEWS 25 Feb 19 Access via Tymnet and SprintNet Eliminated Effective 3/31/02  
NEWS 26 Mar 08 Gene Names now available in BIOSIS

NEWS EXPRESS February 1 CURRENT WINDOWS VERSION IS V6.0d,  
CURRENT MACINTOSH VERSION IS V6.0a(ENG) AND V6.0Ja(JP),  
AND CURRENT DISCOVER FILE IS DATED 05 FEBRUARY 2002

NEWS HOURS STN Operating Hours Plus Help Desk Availability  
NEWS INTER General Internet Information  
NEWS LOGIN Welcome Banner and News Items  
NEWS PHONE Direct Dial and Telecommunication Network Access to STN  
NEWS WWW CAS World Wide Web Site (general information)

Enter NEWS followed by the item number or name to see news on that specific topic.

All use of STN is subject to the provisions of the STN Customer agreement. Please note that this agreement limits use to scientific research. Use for software development or design or implementation of commercial gateways or other similar uses is prohibited and may result in loss of user privileges and other penalties.

FILE 'HOME' ENTERED AT 11:46:00 ON 21 MAR 2002

FILE 'REGISTRY' ENTERED AT 11:46:06 ON 21 MAR 2002  
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.  
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.  
COPYRIGHT (C) 2002 American Chemical Society (ACS)

STRUCTURE FILE UPDATES: 19 MAR 2002 HIGHEST RN 401892-67-9  
DICTIONARY FILE UPDATES: 19 MAR 2002 HIGHEST RN 401892-67-9

TSCA INFORMATION NOW CURRENT THROUGH July 7, 2001

Please note that search-term pricing does apply when conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Calculated physical property data is now available. See HELP PROPERTIES for more information. See STNote 27, Searching Properties in the CAS Registry File, for complete details:  
<http://www.cas.org/ONLINE/STN/STNOTES/stnotes27.pdf>

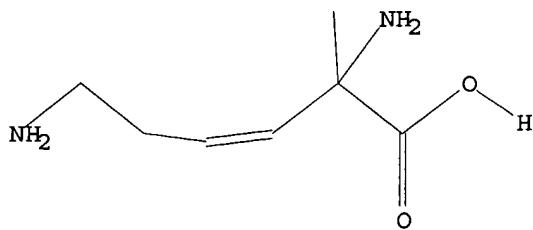
The P indicator for Preparations was not generated for all of the CAS Registry Numbers that were added to the H/Z/CA/CAplus files between 12/27/01 and 1/23/02. Use of the P indicator in online and SDI searches during this period, either directly appended to a CAS Registry Number or by qualifying an L-number with /P, may have yielded incomplete results. As of 1/23/02, the situation has been resolved. Also, note that searches conducted using the PREP role indicator were not affected.

Customers running searches and/or SDIs in the H/Z/CA/CAplus files incorporating CAS Registry Numbers with the P indicator between 12/27/01 and 1/23/02, are encouraged to re-run these strategies. Contact the CAS Help Desk at 1-800-848-6533 in North America or 1-614-447-3698, worldwide, or send an e-mail to [help@cas.org](mailto:help@cas.org) for further assistance or to receive a credit for any duplicate searches.

```
=>  
Uploading 09954988 cl202 last cmpd.str
```

## L1 STRUCTURE UPLOADED

=> d 11  
L1 HAS NO ANSWERS  
L1 . STR



Structure attributes must be viewed using STN Express query preparation.

```
=> search l1 exact full
FULL SEARCH INITIATED 11:46:43 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 13 TO ITERATE
```

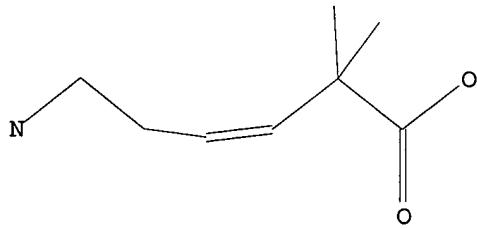
```
100.0% PROCESSED      13 ITERATIONS          0 ANSWERS
SEARCH TIME: 00.00.01
```

```
L2      0 SEA EXA FUL L1
```

```
=>
Uploading 09954988 cl202 last cmpd gnrc.str
```

```
L3      STRUCTURE UPLOADED
```

```
=> d 13
L3 HAS NO ANSWERS
L3      STR
```



Structure attributes must be viewed using STN Express query preparation.

```
=> search l3 sss full
FULL SEARCH INITIATED 11:48:22 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 2987 TO ITERATE
```

```
100.0% PROCESSED      2987 ITERATIONS        0 ANSWERS
SEARCH TIME: 00.00.01
```

```
L4      0 SEA SSS FUL L3
```

```
=>
Uploading 09954988 cl202 last cmpd gnrc.str
```

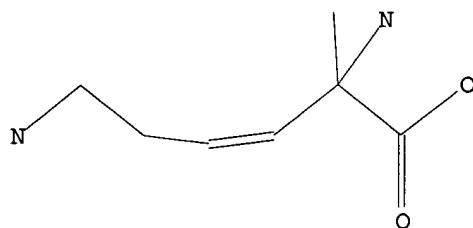
```
L5      STRUCTURE UPLOADED
```

```
=> d 15
```

L5 HAS NO ANSWERS

L5

STR



Structure attributes must be viewed using STN Express query preparation.

```
=> search l5 sss full  
FULL SEARCH INITIATED 11:49:40 FILE 'REGISTRY'  
FULL SCREEN SEARCH COMPLETED - 2201 TO ITERATE
```

100.0% PROCESSED 2201 ITERATIONS  
SEARCH TIME: 00.00.01

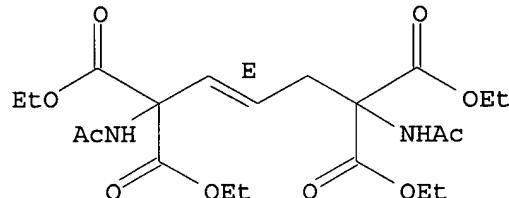
18 ANSWERS

L6 18 SEA SSS FUL L5

=> d scan

L6 18 ANSWERS REGISTRY COPYRIGHT 2002 ACS  
IN 2-Pentene-1,1,5,5-tetracarboxylic acid, 1,5-bis(acetylamino)-, tetraethyl  
ester, (E)- (9CI)  
MF C21 H32 N2 O10

Double bond geometry as shown.



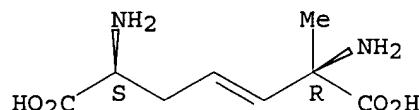
\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):18

L6 18 ANSWERS REGISTRY COPYRIGHT 2002 ACS  
IN 3-Heptenedioic acid, 2,6-diamino-2-methyl-, [R-(R\*,S\*)]- (9CI)  
MF C8 H14 N2 O4

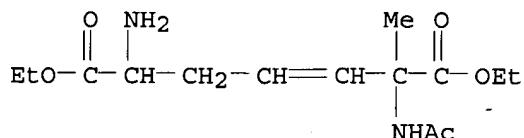
Absolute stereochemistry.

Double bond geometry unknown.



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

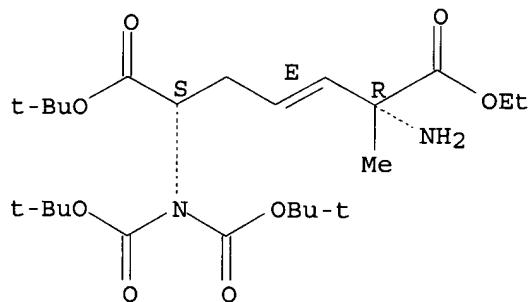
L6 18 ANSWERS REGISTRY COPYRIGHT 2002 ACS  
IN 3-Heptenedioic acid, 2-(acetylamino)-6-amino-2-methyl-, diethyl ester  
(9CI)  
MF C14 H24 N2 O5



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

L6 18 ANSWERS REGISTRY COPYRIGHT 2002 ACS  
IN 3-Heptenedioic acid,  
2-amino-6-[bis[(1,1-dimethylethoxy)carbonyl]amino]-2-  
methyl-, 7-(1,1-dimethylethyl) 1-ethyl ester, [S-[R\*,S\*- (E)]]- (9CI)  
MF C24 H42 N2 O8

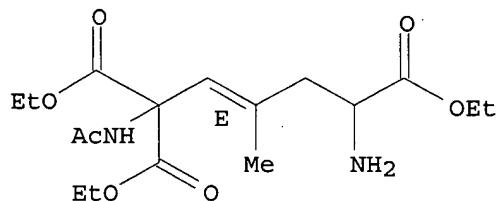
Absolute stereochemistry.  
Double bond geometry as shown.



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

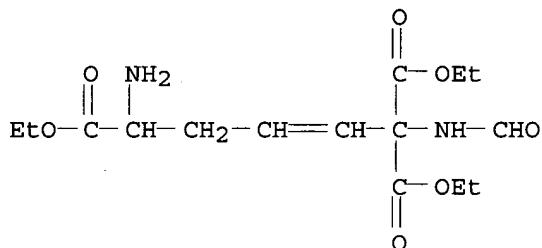
L6 18 ANSWERS REGISTRY COPYRIGHT 2002 ACS  
IN 2-Pentene-1,1,5-tricarboxylic acid, 1-(acetylamino)-5-amino-3-methyl-,  
triethyl ester, monohydrochloride, (E)- (9CI)  
MF C17 H28 N2 O7 . Cl H

Double bond geometry as shown.



● HCl

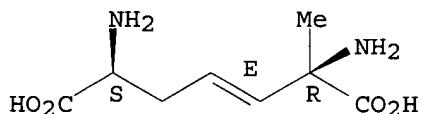
L6 18 ANSWERS REGISTRY COPYRIGHT 2002 ACS  
 IN 2-Pentene-1,1,5-tricarboxylic acid, 5-amino-1-(formylamino)-, triethyl ester (9CI)  
 MF C15 H24 N2 O7



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

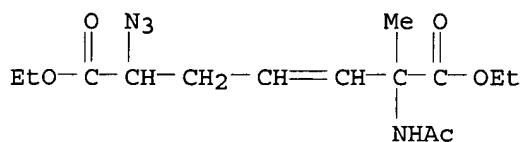
L6 18 ANSWERS REGISTRY COPYRIGHT 2002 ACS  
 IN 3-Heptenedioic acid, 2,6-diamino-2-methyl-, [R-[R\*,S\*- (E)]]- (9CI)  
 MF C8 H14 N2 O4

Absolute stereochemistry.  
 Double bond geometry as shown.

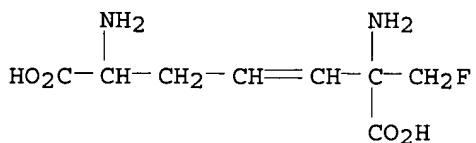


\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

L6 18 ANSWERS REGISTRY COPYRIGHT 2002 ACS  
 IN 3-Heptenedioic acid, 2-(acetylamino)-6-azido-2-methyl-, diethyl ester (9CI)  
 MF C14 H22 N4 O5



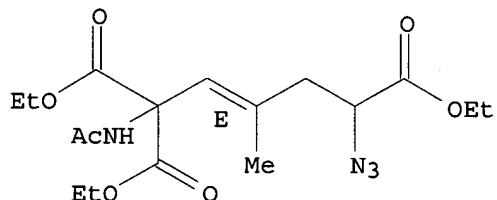
L6 18 ANSWERS REGISTRY COPYRIGHT 2002 ACS  
IN 3-Heptenedioic acid, 2,6-diamino-2-(fluoromethyl)- (9CI)  
MF C8 H13 F N2 O4



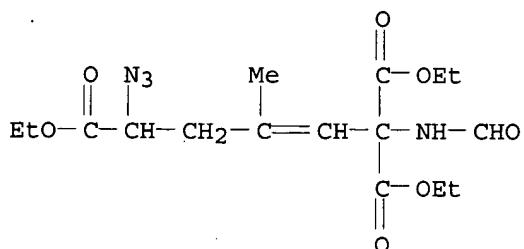
\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

L6 18 ANSWERS REGISTRY COPYRIGHT 2002 ACS  
IN 2-Pentene-1,1,5-tricarboxylic acid, 1-(acetylamino)-5-azido-3-methyl-, triethyl ester, (E)- (9CI)  
MF C17 H26 N4 O7

Double bond geometry as shown.



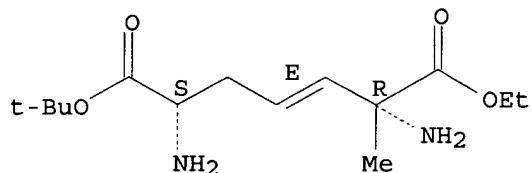
L6 18 ANSWERS REGISTRY COPYRIGHT 2002 ACS  
IN 2-Pentene-1,1,5-tricarboxylic acid, 5-azido-1-(formylamino)-3-methyl-, triethyl ester (9CI)  
MF C16 H24 N4 O7



L6 18 ANSWERS REGISTRY COPYRIGHT 2002 ACS

IN 3-Heptenedioic acid, 2,6-diamino-2-methyl-, 7-(1,1-dimethylethyl) 1-ethyl ester, [S-[R\*,S\*-(E)]]- (9CI)  
MF C14 H26 N2 O4

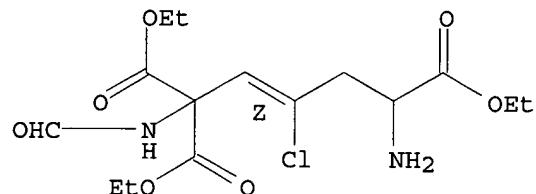
Absolute stereochemistry.  
Double bond geometry as shown.



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

L6 18 ANSWERS REGISTRY COPYRIGHT 2002 ACS  
IN 2-Pentene-1,1,5-tricarboxylic acid, 5-amino-3-chloro-1-(formylamino)-, triethyl ester, (Z)- (9CI)  
MF C15 H23 Cl N2 O7

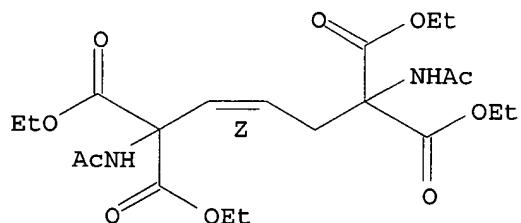
Double bond geometry as shown.



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

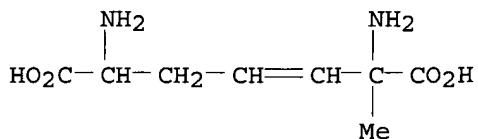
L6 18 ANSWERS REGISTRY COPYRIGHT 2002 ACS  
IN 2-Pentene-1,1,5,5-tetracarboxylic acid, 1,5-bis(acetylamino)-, tetraethyl ester, (Z)- (9CI)  
MF C21 H32 N2 O10

Double bond geometry as shown.



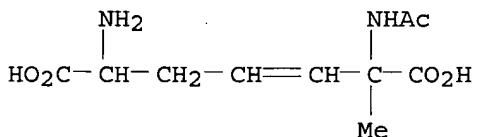
\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

L6 18 ANSWERS REGISTRY COPYRIGHT 2002 ACS  
IN 3-Heptenedioic acid, 2,6-diamino-2-methyl- (9CI)  
MF C8 H14 N2 O4



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

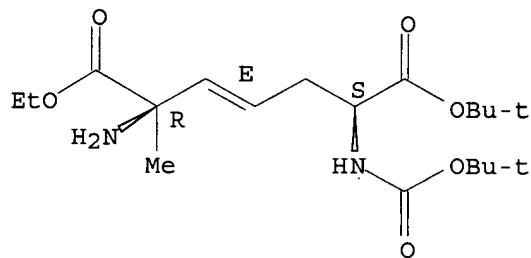
L6 18 ANSWERS REGISTRY COPYRIGHT 2002 ACS  
IN 3-Heptenedioic acid, 2-(acetylamino)-6-amino-2-methyl- (9CI)  
MF C10 H16 N2 O5



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

L6 18 ANSWERS REGISTRY COPYRIGHT 2002 ACS  
IN 3-Heptenedioic acid, 2-amino-6-[(1,1-dimethylethoxy)carbonyl]amino]-2-  
methyl-, 7-(1,1-dimethylethyl) 1-ethyl ester, [S-[R\*,S\*-(E)]]- (9CI)  
MF C19 H34 N2 O6

Absolute stereochemistry.  
Double bond geometry as shown.

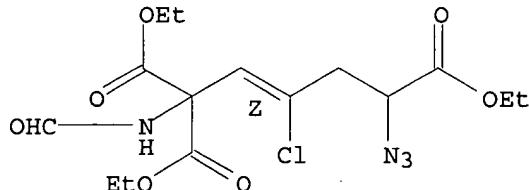


\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

L6 18 ANSWERS REGISTRY COPYRIGHT 2002 ACS  
IN 2-Pentene-1,1,5-tricarboxylic acid, 5-azido-3-chloro-1-(formylamino)-,  
triethyl ester, (Z)- (9CI)

MF C15 H21 Cl N4 O7

Double bond geometry as shown.



ALL ANSWERS HAVE BEEN SCANNED

=> file caplus	COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST		329.68	329.83

FILE 'CAPLUS' ENTERED AT 11:51:21 ON 21 MAR 2002  
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.  
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.  
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FILE COVERS 1907 - 21 Mar 2002 VOL 136 ISS 12  
FILE LAST UPDATED: 19 Mar 2002 (20020319/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

CAS roles have been modified effective December 16, 2001. Please check your SDI profiles to see if they need to be revised. For information on CAS roles, enter HELP ROLES at an arrow prompt or use the CAS Roles thesaurus (/RL field) in this file.

The P indicator for Preparations was not generated for all of the CAS Registry Numbers that were added to the CAS files between 12/27/01 and 1/23/02. As of 1/23/02, the situation has been resolved. Searches and/or SDIs in the H/Z/CA/CAplus files incorporating CAS Registry Numbers with the P indicator executed between 12/27/01 and 1/23/02 may be incomplete. See the NEWS message on this topic for more information.

=> 16  
L7 7 L6

=> d 17 1-7 ti

- L7 ANSWER 1 OF 7 CAPLUS COPYRIGHT 2002 ACS  
TI Purification and characterization of dihydroadipic acid synthase from pea
- L7 ANSWER 2 OF 7 CAPLUS COPYRIGHT 2002 ACS  
TI Stereoselective synthesis of (2S,6S)-2,6-diaminoheptanedioic acid and of unsymmetrical derivatives of meso-2,6-diaminoheptanedioic acid
- L7 ANSWER 3 OF 7 CAPLUS COPYRIGHT 2002 ACS  
TI Synthesis and antimicrobial evaluation of 2,6-diamino-2-fluoromethyl-3-heptenedioic acid, an unsaturated  $\alpha$ -fluoromethyl derivative of diaminopimelic acid
- L7 ANSWER 4 OF 7 CAPLUS COPYRIGHT 2002 ACS  
TI Preparation of (Z)-3,4-didehydronorvaline, (E)- and (Z)-3,4-didehydroornithine and (E)- and (Z)-3,4-didehydro-2,6-diaminopimelic acid
- L7 ANSWER 5 OF 7 CAPLUS COPYRIGHT 2002 ACS  
TI Preparation of (glutamylamino)alkanedioates as autoimmune medicine and pharmaceutical compositions containing them
- L7 ANSWER 6 OF 7 CAPLUS COPYRIGHT 2002 ACS  
TI Diaminopimelic acid derivatives
- L7 ANSWER 7 OF 7 CAPLUS COPYRIGHT 2002 ACS  
TI The lysine pathway as a target for a new genera of synthetic antibacterial antibiotics?

=> logoff hold  
COST IN U.S. DOLLARS  
FULL ESTIMATED COST

	SINCE FILE	TOTAL
	ENTRY	SESSION
FULL ESTIMATED COST	5.34	335.17

SESSION WILL BE HELD FOR 60 MINUTES  
STN INTERNATIONAL SESSION SUSPENDED AT 11:57:29 ON 21 MAR 2002  
Connection closed by remote host

Connecting via Winsock to STN

Trying 3106016892...Open

Welcome to STN International! Enter x:x  
LOGINID:ssspta1623paz  
PASSWORD:  
\* \* \* \* \* RECONNECTED TO STN INTERNATIONAL \* \* \* \* \*  
SESSION RESUMED IN FILE 'CAPLUS' AT 12:12:53 ON 21 MAR 2002  
FILE 'CAPLUS' ENTERED AT 12:12:53 ON 21 MAR 2002  
COPYRIGHT (C) 2002 AMERICAN CHEMICAL SOCIETY (ACS)

COST IN U.S. DOLLARS  
FULL ESTIMATED COST

	SINCE FILE	TOTAL
	ENTRY	SESSION
FULL ESTIMATED COST	5.34	335.17

=> save temp 17 cl202srch/  
CL202SRCH/ IS NOT A VALID SAVED NAME  
Enter the name you wish to use for the saved query,

answer set, or L-number list. The name must:

1. Begin with a letter,
2. Have 1-12 characters,
3. Contain only letters (A-Z) and numbers (0-9),
4. End with /Q for a query (search profile, structure, or screen set), /A for an answer set, or /L for an L-number list.
5. Not already be in use as a saved name,
6. Not be END, SAV, SAVE, SAVED
7. Not have the form of an L-number (Lnnn).

ENTER NAME OR (END) :end

=> save temp l7 cl202srch/a  
ANSWER SET L7 HAS BEEN SAVED AS 'CL202SRCH/A'

=> logoff hold

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	5.68	335.51

SESSION WILL BE HELD FOR 60 MINUTES  
STN INTERNATIONAL SESSION SUSPENDED AT 12:13:40 ON 21 MAR 2002

L Number	Hits	Search Text	DB	Time stamp
1	0	9313055.URPN.	USPAT; EPO; JPO; DERWENT	2002/03/21 13:12
2	337	562/560.ccls.	USPAT; EPO; JPO; DERWENT	2002/03/21 13:12
3	389	562/561.ccls.	USPAT; EPO; JPO; DERWENT	2002/03/21 13:12
4	244	562/564.ccls.	USPAT; EPO; JPO; DERWENT	2002/03/21 13:12
5	862	562/560.ccls. or 562/561.ccls. or 562/564.ccls.	USPAT; EPO; JPO; DERWENT	2002/03/21 13:12
6	9	nitric adj oxide	USPAT; EPO; JPO; DERWENT	2002/03/21 13:12
7	7939	nitric adj oxide	USPAT; EPO; JPO; DERWENT	2002/03/21 13:12
8	0	methylserinate	USPAT; EPO; JPO; DERWENT	2002/03/21 13:12
9	173	heptynoic	USPAT; EPO; JPO; DERWENT	2002/03/21 13:12
10	1461	heptenoic	USPAT; EPO; JPO; DERWENT	2002/03/21 13:12
11	3	9525717.pn.	USPAT; EPO; JPO; DERWENT	2002/03/21 13:12
12	2	9313055.pn.	USPAT; EPO; JPO; DERWENT	2002/03/21 13:12
13	2	5863931.URPN.	USPAT; EPO; JPO; DERWENT	2002/03/21 13:12
14	3	9533534.pn.	USPAT; EPO; JPO; DERWENT	2002/03/21 13:12
15	2	9534534.pn.	USPAT; EPO; JPO; DERWENT	2002/03/21 13:12
16	5	9104024.pn.	USPAT; EPO; JPO; DERWENT	2002/03/21 13:12
17	3	5798387.pn.	USPAT; EPO; JPO; DERWENT	2002/03/21 13:12
18	24	(nitric adj oxide) and (562/560.ccls. or 562/561.ccls. or 562/564.ccls.)	USPAT; EPO; JPO; DERWENT	2002/03/21 13:12
19	4	3402990.pn.	USPAT; EPO; JPO; DERWENT	2002/03/21 13:12
20	3	5132453.pn.	USPAT; EPO; JPO; DERWENT	2002/03/21 13:12
21	2	(562/560.ccls. or 562/561.ccls. or 562/564.ccls.) and heptynoic	USPAT; EPO; JPO; DERWENT	2002/03/21 13:12
22	1	9324126.pn.	USPAT; EPO; JPO; DERWENT	2002/03/21 13:12
23	7	heptenoic and (562/560.ccls. or 562/561.ccls. or 562/564.ccls.)	USPAT; EPO; JPO; DERWENT	2002/03/21 13:12

24	3	5854234.pn.	USPAT; EPO; JPO; DERWENT USPAT; EPO; JPO; DERWENT	2002/03/21 13:12
25	2	6046211.pn.		2002/03/21 13:12

	Type	L #	Hits	Search Text	DBs	Time Stamp	Comments	Error Definition
1	BRS	L1	0	9313055.URPN.	USPAT ; EPO; JPO; DERWE NT	2002/03/21 13:12		
2	BRS	L2	337	562/560.ccls.	USPAT ; EPO; JPO; DERWE NT	2002/03/21 13:12		
3	BRS	L3	389	562/561.ccls.	USPAT ; EPO; JPO; DERWE NT	2002/03/21 13:12		
4	BRS	L4	244	562/564.ccls.	USPAT ; EPO; JPO; DERWE NT	2002/03/21 13:12		
5	BRS	L5	862	562/560.ccls. or 562/561.ccls. or 562/564.ccls.	USPAT ; EPO; JPO; DERWE NT	2002/03/21 13:12		
6	BRS	L6	9	nitic adj oxide	USPAT ; EPO; JPO; DERWE NT	2002/03/21 13:12		
7	BRS	L7	7939	nitric adj oxide	USPAT ; EPO; JPO; DERWE NT	2002/03/21 13:12		
8	BRS	L8	0	methylserinate	USPAT ; EPO; JPO; DERWE NT	2002/03/21 13:12		
9	BRS	L9	173	heptynoic	USPAT ; EPO; JPO; DERWE NT	2002/03/21 13:12		
10	BRS	L10	1461	heptenoic	USPAT ; EPO; JPO; DERWE NT	2002/03/21 13:12		

	Errors
1	0
2	0
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	Type	L #	Hits	Search Text	DBs	Time Stamp	Comments	Error Definition
11	BRS	L11	3	9525717.pn.	USPAT ; EPO; JPO; DERWE NT	2002/03/21 13:12		
12	BRS	L12	2	9313055.pn.	USPAT ; EPO; JPO; DERWE NT	2002/03/21 13:12		
13	BRS	L13	2	5863931.URPN.	USPAT ; EPO; JPO; DERWE NT	2002/03/21 13:12		
14	BRS	L14	3	9533534.pn.	USPAT ; EPO; JPO; DERWE NT	2002/03/21 13:12		
15	BRS	L15	2	9534534.pn.	USPAT ; EPO; JPO; DERWE NT	2002/03/21 13:12		
16	BRS	L16	5	9104024.pn.	USPAT ; EPO; JPO; DERWE NT	2002/03/21 13:12		
17	BRS	L17	3	5798387.pn.	USPAT ; EPO; JPO; DERWE NT	2002/03/21 13:12		
18	BRS	L18	24	(nitric adj oxide) and (562/560.ccls. or 562/561.ccls. or 562/564.ccls.)	USPAT ; EPO; JPO; DERWE NT	2002/03/21 13:12		
19	BRS	L19	4	3402990.pn.	USPAT ; EPO; JPO; DERWE NT	2002/03/21 13:12		
20	BRS	L20	3	5132453.pn.	USPAT ; EPO; JPO; DERWE NT	2002/03/21 13:12		

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	Type	L #	Hits	Search Text	DBs	Time Stamp	Comments	Error Definition
21	BRS	L21	2	(562/560.ccls. or 562/561.ccls. or 562/564.ccls.) and heptenoic	USPAT ; EPO; JPO; DERWE NT	2002/03/21 13:12		
22	BRS	L22	1	9324126.pn.	USPAT ; EPO; JPO; DERWE NT	2002/03/21 13:12		
23	BRS	L23	7	heptenoic and (562/560.ccls. or 562/561.ccls. or 562/564.ccls.)	USPAT ; EPO; JPO; DERWE NT	2002/03/21 13:12		
24	BRS	L24	3	5854234.pn.	USPAT ; EPO; JPO; DERWE NT	2002/03/21 13:12		
25	BRS	L25	2	6046211.pn.	USPAT ; EPO; JPO; DERWE NT	2002/03/21 13:12		

	Err ors
21	0
22	0
23	0
24	0
25	0